



UNITED STATES ENVIRONMENTAL PROTECTION AGENCY

REGION 7  
25 FUNSTON ROAD  
KANSAS CITY, KANSAS 66115

RECEIVED

SEP 24 1991

PRMT SECTION

SEP 20 1991

MEMORANDUM

SUBJECT: Transmittal of Analytical Data: Monsanto-Queeny Plant  
St. Louis, Missouri (ADF43)

FROM: Robert B. Dona *RBDona*  
Environmental Engineer, QADE/EDSB/ENSV

TO: Pat Nichols  
Work Assignment Manager, PRMT/RCRA/WSTM

THRU: Jeffrey A. Wandtke *JAW*  
Regional QA Officer, ENSV

I am transmitting the attached Analysis Request Report for the RCRA groundwater sampling at the Monsanto-Queeny Plant, St. Louis, Missouri, by Steve Freeman of the Jacobs Engineering Group Inc. on June 4-7, 1991. This Analysis Request Report contains only the data from the groundwater samples collected as split samples with the facility's contractor and the field QC samples. The data from the analysis of the laboratory quality control are available at your request.

I am also including copies of the field sheets, the chain-of-custody record and ESAT case narratives and data reviews. If you have any questions, please call me at 551-5182.

Attachments



R00105659  
RCRA RECORDS CENTER

ACCH09

## ANALYSIS REQUEST REPORT

VALIDATED DATA

FOR ACTIVITY: ADF43

DONA, B.

09/18/91 17:19:19

ALL REAL SAMPLES AND FIELD Q.C.

## \* FINAL REPORT

FY: 91 ACTIVITY: ADF43 DESCRIPTION: MONSANTO-QUEENY LOCATION: ST. LOUIS MISSOURI

STATUS: ACTIVE TYPE: SAMPLING - IN HOUSE ANALYSIS PROJECT: A60

LABO DUE DATE IS 7/10/91. REPORT DUE DATE IS 7/28/91.

INSPECTION DATE: 6/ 7/91 ALL SAMPLES RECEIVED DATE: 06/10/91

ALL DATA APPROVED BY LABO DATE: 09/10/91 FINAL REPORT TRANSMITTED DATE: 00/00/00

EXPECTED LABO TURNAROUND TIME IS 30 DAYS EXPECTED REPORT TURNAROUND TIME IS 51 DAYS

ACTUAL LABO TURNAROUND TIME IS 92 DAYS ACTUAL REPORT TURNAROUND TIME IS 0 DAYS

SAMP. NO.	QCC	M	DESCRIPTION	SAMPLE STATUS	# CONT.	CITY	STATE	AIRS/ STORET LOC NO	BEG. DATE	BEG. TIME	END. DATE	END. TIME
001		W	MONSANTO-QUEENY-MW-11	1	8	ST. LOUIS	MISSOURI		06/04/91	10:30	/ /	:
002		W	MONSANTO-QUEENY-TW-1	1	8	ST. LOUIS	MISSOURI		06/04/91	17:20	/ /	:
004		W	MONSANTO-QUEENY-QS-1	1	8	ST. LOUIS	MISSOURI		06/04/91	:	/ /	:
005		W	MONSANTO QUEENY-HW-1B	1	8	ST. LOUIS	MISSOURI		06/05/91	14:30	06/05/91	15:20
006		W	MONSANTO QUEENY-VW-2B	1	8	ST. LOUIS	MISSOURI		06/06/91	09:00	06/06/91	09:30
007		W	MONSANTO QUEENY-MW-8A	1	8	ST. LOUIS	MISSOURI		06/07/91	09:15	06/07/91	09:45
008	F	W	MONSANTO QUEENY-RINSATE BLANK	1	3	ST. LOUIS	MISSOURI		06/07/91	08:20	/ /	:
009	F	W	MONSANTO-QUEENY=FIELD BLANK	1	1	ST. LOUIS	MISSOURI		06/04/91	17:10	/ /	:

## TABLE OF CODES

## VALIDATED DATA

SAMP. NO. = SAMPLE IDENTIFICATION NUMBER  
 QCC = QUALITY CONTROL SAMPLE/AUDIT CODE  
 M = MEDIA OF SAMPLE (A=AIR, T=TISSUE, H=HAZARDOUS MATERIAL, S=SEDIMENT/SOIL, W=WATER)

AIRS/STORET LOC. NO. = A SAMPLING SITE LOCATION IDENTIFICATION NUMBER

BEG. DATE = THE DATE SAMPLING WAS STARTED

BEG. TIME = THE TIME SAMPLING WAS STARTED

END. DATE = THE DATE SAMPLING WAS ENDED

END. TIME = THE TIME SAMPLING WAS STOPPED

A = RESERVED

B = RESERVED

PES = PESTICIDES BY CONTRACT

= DIOXINS/FURANS BY EPA

E = EXPLOSIVES BY CONTRACT

FLD = FIELD MEASUREMENTS BY EPA

G = MINERALS & DISSOLVED MATERIALS BY EPA

HER = HERBICIDES BY EPA

I = ION CHROMATOGRAPHY ANALYSES BY EPA

MC = METALS BY CONTRACT

BNC = BASE NEUTRALS BY CONTRACT

L = FISH PHYSICAL DATA BY EPA

MET = METALS BY EPA

N = FISH TISSUE PARAMETERS BY EPA

VC = VOLATILES BY CONTRACT

P = PESTICIDES BY EPA

Q = FLASH POINT ANALYSES BY EPA

R = RESERVED

BN = SEMIVOLATILE BY EPA

T = CYANIDE PHENOL BY EPA

U = RESERVED

VOA = VOLATILE ORGANICS BY EPA

HC = HERBICIDES BY CONTRACT

X = RESERVED

Y = RESERVED

TRK = ACTIVITY TRACKING PARAMETERS BY EPA

## STORET DETECTION IDENTIFIERS

BLANK = NO REMARKS

J = DATA REPORTED BUT NOT VALID BY APPROVED QC PROCEDURES

I = INVALID SAMPLE/DATA - VALUE NOT REPORTED

U = LESS THAN (MEASUREMENT DETECTION LIMIT)

M = DETECTED BUT BELOW THE LEVEL FOR ACCURATE QUANTIFICATION

O = PARAMETER NOT ANALYZED

## CONTRACTOR/ IN HOUSE / FIELD MEDIA GROUPS

FIELD = \* \* \* = AF, HF, SF, TF, WF, ZZ

CONTRACTOR = \* \* = HA, HC, HJ, HK, HO, SC, SJ, SK, SO, SW, TC, TJ, TK, TO, TW, WA, WC, WE, WJ, WK, WO, WW

IN HOUSE = \* = ALL OTHERS

## QUALITY CONTROL AUDIT CODES

A = TRUE VALUE FOR CALIBRATION STANDARD

B = CONCENTRATION RESULTING FROM DUPLICATE LAB SPIKE

C = MEASURED VALUE FOR CALIBRATION STANDARD

D = MEASURED VALUE FOR FIELD DUPLICATE

F = MEASURED VALUE FOR FIELD BLANK

G = MEASURED VALUE FOR METHOD STANDARD

H = TRUE VALUE FOR METHOD STANDARD

K = CONCENTRATION RESULTING FROM DUPLICATE FIELD SPIKE

L = MEASURED VALUE FOR LAB DUPLICATE

M = MEASURED VALUE FOR LAB BLANK

N = MEASURED VALUE FOR DUPLICATE FIELD SPIKE

P = MEASURED VALUE FOR PERFORMANCE STANDARD

R = CONCENTRATION RESULTING FROM LAB SPIKE

S = MEASURED VALUE FOR LAB SPIKE

T = TRUE VALUE OF PERFORMANCE STANDARD

W = MEASURED VALUE FOR DUPLICATE LAB SPIKE

Y = MEASURED VALUE FOR FIELD SPIKE

Z = CONCENTRATION RESULTING FROM FIELD SPIKE

## MEDIA CODES

A = AIR

T = BIOLOGICAL (PLANT & ANIMAL) TISSUE

H = HAZARDOUS MATERIALS/MAN MADE PRODUCTS

S = SEDIMENT, SLUDGE & SOIL

W = WATER

## UNITS

NA = NOT APPLICABLE

PG = PICOGRAMS (1 X 10<sup>-12</sup> GRAMS)

NG = NANOGRAMS (1 X 10<sup>-9</sup> GRAMS)

UG = MICROGRAMS (1 X 10<sup>-6</sup> GRAMS)

MG = MILLIGRAMS (1 X 10<sup>-3</sup> GRAMS)

M3 = METER CUBED

MPH = MILES PER HOUR

SCM = STANDARD (1 ATM, 25 C) CUBIC METER

KG = KILOGRAM

L = LITER

C = CENTIGRADE DEGREES

SU = STANDARD (PH) UNITS

# = NUMBER

LB = POUNDS

IN = INCHES

M/F = MALE/FEMALE

M2 = SQUARE METER

I.D. = SPECIES IDENTIFICATION

GPM = GALLONS PER MINUTE

CFS = CUBIC FEET PER SECOND

MGD = MILLION GALLONS PER DAY

1000G = FLOW, 1000 GALLONS PER COMPOSITE

UMHOS = CONDUCTIVITY UNITS (1/OHMS)

NTU = TURBIDITY UNITS

PC/L = PICO (1 X 10<sup>-12</sup>) CURRIES PER LITER

MV = MILLIVOLT

SQ FT = SQUARE FEET

P/CM2 = PICOGRAMS PER SQ. CENTIMETER

U/CM2 = MICROGRAMS PER SQ. CENTIMETER

## ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 1-ADF43

VALIDATED DATA

COMPOUND	UNITS	001	002	004	005	006
WD02 2,3,7,8-TCDD IN WATER (PG/ML)	PG/ML	0.000174 U	0.00035 U	0.000176 U	0.00057 U	0.00047 U
WF01 WATER TEMP	'C	76.0	75.8	67.01	68.7	65.5
WF05 PH, FIELD	SU	6.93	6.75	8.06	7.88	6.33
WF10 CONDUCTIVITY (FIELD)	UMHOS	1210	2550	3730	1953	776
WH01 2,4-D	UG/L	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
WH02 2,4,5-TP (SILVEX)	UG/L	0.2 U	0.2 U	11 U	0.2 U	0.2 U
WH03 2,4,5-T	UG/L	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
WM01 SILVER BY ICAP	UG/L	10 U	10 U	10 U	10 U	10 U
WM02 ALUMINUM, TOTAL, BY ICAP	UG/L	681	258	5990	4990	5850
WM03 ARSENIC, TOTAL, BY ICAP	UG/L	50 U	50 U	50 U	50 U	50 U
WM04 BARIUM, TOTAL, BY ICAP	UG/L	285	266	2030	219	175
WM05 BERYLLIUM, TOTAL, BY ICAP	UG/L	2 U	2 U	2 U	2 U	2 U
WM06 CADMIUM, TOTAL, BY ICAP	UG/L	5 U	5 U	5 U	5 U	5 U
WM07 COBALT, TOTAL, BY ICAP	UG/L	10 U	10 U	10 U	10 U	10 U
WM08 CHROMIUM, TOTAL, BY ICAP	UG/L	10 U	10 U	14.6	18.2	10 U
WM09 COPPER, TOTAL, BY ICAP	UG/L	10 U	10 U	34.2	21.5	17.9
WM10 IRON BY ICAP	UG/L	10200	31300	17000	9300	27900
WM11 MANGANESE BY ICAP	UG/L	385	2280	449	371	1500
WM12 MOLYBDENUM BY ICAP	UG/L	NA 0	NA 0	NA 0	NA 0	NA 0
WM13 NICKEL BY ICAP	UG/L	20 U	20 U	20 U	21.8	20 U
WM14 LEAD BY ICAP	UG/L	50 U	50 U	90	50 U	50 U
WM15 ANTIMONY, TOTAL, BY ICAP	UG/L	50 U	50 U	50 U	50 U	50 U
WM16 SELENIUM BY ICAP	UG/L	54.7	64.3	50 U	50 U	50 U
WM17 TITANIUM BY ICAP	UG/L	NA 0	NA 0	NA 0	NA 0	NA 0
WM18 THALLIUM BY ICAP	UG/L	300 U	300 U	300 U	300 U	300 U
WM19 VANADIUM BY ICAP	UG/L	10 U	10 U	19.1	15.1	15.4

## ANALYSIS REQUEST DETAIL REPORT

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VALIDATED DATA

COMPOUND	UNITS	001	002	004	005	006
WM20 ZINC BY ICAP	UG/L	37.2	122	153	89	42.9
WM21 CALCIUM, TOTAL BY ICAP	MG/L	179	262	59.4	52.1	191
WM22 MAGNESIUM, TOTAL BY ICAP	MG/L	14.7	63.8	20	6.99	40.7
WM23 SODIUM, TOTAL BY ICAP	MG/L	33.5	172	565	149	76.1
WM24 POTASSIUM, TOTAL BY ICAP	MG/L	5.83	36.2	63.8	343	23.5
WM34 MERCURY BY COLD VAPOR AA	UG/L	0.2 U	0.218	0.598	0.2 U	0.2 U
WP01 BHC, ALPHA, BY GC/EC	UG/L	5.0 U	0.05 U	0.050 U	0.050 U	0.050 U
WP02 BHC, BETA, BY GC/EC	UG/L	5.0 U	0.05 U	0.050 U	0.050 U	0.050 U
WP03 BHC, DELTA	UG/L	5.0 U	0.05 U	0.050 U	0.050 U	0.050 U
WP04 GAMMA-BHC(LINDANE), BY GC/EC	UG/L	5.0 U	0.05 U	0.050 U	0.050 U	0.050 U
WP05 ALDRIN, BY GC/EC	UG/L	5.0 U	0.05 U	0.050 U	0.050 U	0.050 U
WP06 DIELDRIN, BY GC/EC	UG/L	10 U	0.10 U	0.10 U	0.10 U	0.10 U
WP07 ENDOSULFAN I, BY GC/EC	UG/L	5.0 U	0.05 U	0.050 U	0.050 U	0.050 U
WP08 ENDOSULFAN II, BY GC/EC	UG/L	10 U	0.10 U	0.10 U	0.10 U	0.10 U
WP09 ENDOSULFAN SULFATE, BY GC/EC	UG/L	10 U	0.10 U	0.10 U	0.10 U	0.10 U
WP10 ENDRIN, BY GC/EC	UG/L	10 U	0.10 U	0.10 U	0.10 U	0.10 U
WP11 ENDRIN ALDEHYDE, BY GC/EC	UG/L	NA O	NA O	NA O	NA O	NA O
WP13 4,4'-DDE	UG/L	10 U	0.10 U	0.10 U	0.10 U	0.10 U
WP14 4,4'-DDD	UG/L	10 U	0.10 U	0.10 U	0.10 U	0.10 U
WP15 4,4'-DDT	UG/L	10 U	0.10 U	0.10 U	0.10 U	0.10 U
WP16 TOXAPHENE	UG/L	100 U	1.0 U	1.0 U	1.0 U	1.0 U
WP17 PCB-1016	UG/L	50 U	0.50 U	0.50 U	0.50 U	0.50 U
WP18 PCB-1221	UG/L	50 U	0.50 U	0.50 U	0.50 U	0.50 U
WP19 PCB-1232	UG/L	50 U	0.50 U	0.50 U	0.50 U	0.50 U
WP20 PCB-1242	UG/L	50 U	0.50 U	0.50 U	0.50 U	0.50 U
WP21 PCB-1248	UG/L	50 U	0.50 U	0.50 U	0.50 U	0.50 U

## ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 1-ADF43

VALIDATED DATA

COMPOUND	UNITS	001	002	004	005	006					
WP22 PCB-1254	UG/L	100	U	1.0	U	1.0	U	1.0	U	1.0	U
WP23 PCB-1260	UG/L	100	U	1.0	U	1.0	U	1.0	U	1.0	U
WP24 CHLORDANE, TECHNICAL, BY GC/EC	UG/L	NA	O	NA	O	NA	O	NA	O	NA	O
WP25 HEPTACHLOR, BY GC/EC	UG/L	5.0	U	0.050	U	0.050	U	0.050	U	0.050	U
WP26 HEPTACHLOR EPOXIDE, BY GC/EC	UG/L	5.0	U	0.050	U	0.050	U	0.050	U	0.050	U
WS01 PHENOL	UG/L	100	U	10	U	10	U	10	U	10	U
WS03 ETHER, BIS(2-CHLOROETHYL), BY GC/MS	UG/L	100	U	190		10	U	10	U	10	U
WS04 2-CHLOROPHENOL	UG/L	100	U	10	U	10	U	10	U	10	U
WS05 1,3-DICHLOROBENZENE	UG/L	100	U	10	U	10	U	10	U	10	U
WS06 1,4-DICHLOROBENZENE	UG/L	100	U	10	U	10	U	10	U	10	U
WS07 BENZYL ALCOHOL	UG/L	100	U	10	U	10	U	10	U	10	U
WS08 1,2-DICHLOROBENZENE	UG/L	100	U	18		10	U	10	U	10	U
WS09 2-METHYLPHENOL (O-CRESOL)	UG/L	100	U	10	U	10	U	10	U	10	U
WS10 ETHER, BIS(2-CHLOROISOPROPYL), BY GC/MS	UG/L	100	U	10	U	10	U	10	U	10	U
WS11 4-METHYLPHENOL (P-CRESOL)	UG/L	100	U	10	U	10	U	10	U	10	U
WS12 N-NITROSO-DIPROPYLAMINE	UG/L	100	U	10	U	10	U	10	U	10	U
WS13 HEXACHLOROETHANE	UG/L	100	U	10	U	10	U	10	U	10	U
WS14 NITROBENZENE	UG/L	100	U	10	U	10	U	10	U	10	U
WS15 ISOPHORONE	UG/L	100	U	10	U	10	U	10	U	10	U
WS16 2-NITROPHENOL	UG/L	100	U	10	U	10	U	10	U	10	U
WS17 2,4-DIMETHYLPHENOL	UG/L	100	U	10	U	10	U	10	U	10	U
WS18 BENZOIC ACID, BY GC/MS	UG/L	500	U	50	U	50	U	50	U	50	U
WS19 METHANE, BIS(2-CHLOROETHOXY), BY GC/MS	UG/L	100	U	10	U	10	U	10	U	10	U
WS20 2,4-DICHLOROPHENOL	UG/L	100	U	10	U	10	U	10	U	10	U
WS21 1,2,4-TRICHLOROBENZENE	UG/L	100	U	10	U	10	U	10	U	10	U
WS22 NAPHTHALENE	UG/L	100	U	10	U	10	U	10	U	10	U

## ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 1-ADF43

VALIDATED DATA

COMPOUND	UNITS	001	002	004	005	006
WS23 4-CHLOROANILINE	UG/L	100 U	10 U	10 U	10 U	10 U
WS24 HEXACHLOROBUTADIENE, BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS25 4-CHLORO-3-METHYLPHENOL	UG/L	100 U	10 U	10 U	10 U	10 U
WS26 2-METHYLNAPHTHALENE	UG/L	100 U	10 U	10 U	10	10 U
WS27 HEXACHLOROCYCLOPENTADIENE, BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS28 2,4,6-TRICHLOROPHENOL	UG/L	100 U	10 U	10 U	10 U	50 U
WS29 2,4,5-TRICHLOROPHENOL	UG/L	500 U	50 U	50 U	50 U	10 U
WS30 2-CHLORONAPHTHALENE	UG/L	100 U	10 U	10 U	10 U	50 U
WS31 2-NITROANILINE (ORTHO NITROANILINE)	UG/L	500 U	50 U	50 U	50 U	10 U
WS32 PHTHALATE, DIMETHYL, BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS33 ACENAPHTHYLENE, BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS34 3-NITROANILINE	UG/L	500 U	50 U	50 U	50 U	50 U
WS35 ACENAPHTHENE, BY GC/MS	UG/L	100 U	10 U	10	10 U	10 U
WS36 2,4-DINITROPHENOL	UG/L	500 U	50 U	50 U	50 U	50 U
WS37 4-NITROPHENOL	UG/L	500 U	50 U	50 U	50 U	50 U
WS38 DIBENZOFURAN	UG/L	100 U	10 U	10 U	10 U	10 U
WS39 2,4-DINITROTOLUENE	UG/L	100 U	10 U	10 U	10 U	10 U
WS40 2,6-DINITROTOLUENE	UG/L	100 U	10 U	10 U	10 U	10 U
WS41 PHTHALATE, DIETHYL, BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS42 4-CHLOROPHENYL PHENYL ETHER	UG/L	100 U	10 U	10 U	10 U	10 U
WS43 FLUORENE, BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS44 4-NITROANILINE	UG/L	500 U	NA I	NA I	NA I	NA I
WS45 4,6-DINITRO-2-METHYLPHENOL	UG/L	500 U	50 U	50 U	50 U	50 U
WS46 N-NITROSODIPHENYLAMINE	UG/L	100 U	10 U	10 U	10 U	10 U
WS47 4-BROMOPHENYL PHENYL ETHER	UG/L	100 U	10 U	10 U	10 U	10 U
WS48 HEXACHLOROBENZENE, BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U

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VALIDATED DATA

COMPOUND	UNITS	001	002	004	005	006
WV10 1,2-DICHLOROETHENE, TOTAL	UG/L	50. U	50.	5.0 U	440.	1900.
WV11 CHLOROFORM, BY GC/MS	UG/L	50. U	5.0 U	5.0 U	5.0 U	5.0 U
WV12 1,2-DICHLOROETHANE	UG/L	50. U	5.0 U	5.0 U	5.0 U	5.0 U
WV13 1,1,1-TRICHLOROETHANE	UG/L	50. U	5.0 U	5.0 U	6.2	5.0 U
WV14 CARBON TETRACHLORIDE, BY GC/MS	UG/L	50. U	5.0 U	5.0 U	5.0 U	5.0 U
WV15 BROMODICHLOROMETHANE, BY GC/MS	UG/L	50. U	5.0U	5.0 U	5.0 U	5.0 U
WV16 1,2-DICHLOROPROPANE	UG/L	50. U	5.0U	5.0 U	5.0 U	5.0 U
WV17 BENZENE, BY GC/MS	UG/L	50. U	60.	5.0 U	5.0 U	5.0 U
WV19 TRICHLOROETHENE	UG/L	50. U	5.0U	5.0 U	1200.	5.0 U
WV20 DICHLOROPROPENE, CIS-1,3-, BY GC/MS	UG/L	50. U	5.0U	5.0 U	5.0 U	5.0 U
WV21 DIBROMOCHLOROMETHANE, BY GC/MS	UG/L	50. U	5.0U	5.0 U	5.0 U	5.0 U
WV22 1,1,2-TRICHLOROETHANE	UG/L	50. U	5.0U	5.0 U	5.0 U	5.0 U
WV24 BROMOFORM, BY GC/MS	UG/L	50. U	5.0U	5.0 U	5.0 U	5.0 U
WV25 TETRACHLOROETHENE	UG/L	50. U	5.0U	5.0 U	5.0 U	5.0 U
WV26 TOLUENE	UG/L	50. U	5.0U	5.0 U	5.0 U	5.0 U
WV27 1,1,2,2-TETRACHLOROETHANE	UG/L	50. U	5.0U	5.0 U	5.0 U	5.0 U
WV28 CHLOROBENZENE, BY GC/MS	UG/L	66000.	35.	6.1	5.0 U	5.0 U
WV29 ETHYL BENZENE, BY GC/MS	UG/L	50. U	5.0U	5.0 U	5.0 U	5.0 U
WV30 ACETONE, BY GC/MS	UG/L	100. U	10. U	10. U	10. U	10. U
WV31 CARBON DISULFIDE, BY GC/MS	UG/L	50. U	5.0U	5.0 U	5.0 U	5.0 U
WV32 2-BUTANONE	UG/L	990.	10. U	10. U	10. U	10. U
WV33 VINYL ACETATE	UG/L	100. U	10. U	10. U	10. U	10. U
WV34 2-HEXANONE	UG/L	100. U	10. U	10. U	10. U	10. U
WV35 4-METHYL-2-PENTANONE	UG/L	100. U	10. U	10. U	10. U	10. U
WV36 STYRENE	UG/L	50. U	5.0U	5.0U	5.0U	5.0U
WV37 XYLENES, TOTAL	UG/L	50. U	5.0U	5.0U	5.0U	5.0U



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COMPOUND	UNITS	001	002	004	005	006
WS49 PENTACHLOROPHENOL	UG/L	500 U	50 U	50 U	50 U	50 U
WS50 PHENANTHRENE	UG/L	100 U	10 U	20 U	10 U	10 U
WS51 ANTHRACENE, BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS52 PHTHALATE, DI-N-BUTYL-, BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS53 FLUORANTHENE, BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS54 PYRENE	UG/L	100 U	10 U	10 U	10 U	10 U
WS55 PHTHALATE, BUTYL BENZYL	UG/L	100 U	10 U	10 U	10 U	10 U
WS56 3,3'-DICHLOROBENZIDINE	UG/L	200 U	20 U	20 U	20 U	20 U
WS57 ANTHRACENE, BENZO(A), BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS58 PHTHALATE, BIS(2-ETHYLHEXYL), BY GC/MS	UG/L	100 U	14 U	16 U	10 U	21 U
WS59 CHRYSENE, BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS60 PHTHALATE, DI-N-OCTYL-, BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS61 FLUORANTHENE, BENZO(B), BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS62 FLUORANTHENE, BENZO(K), BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS63 PYRENE, BENZO(A), BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS64 INDENO(1,2,3-CD)PYRENE	UG/L	100 U	10 U	10 U	10 U	10 U
WS65 ANTHRACENE, DIBENZO(A,H), BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS66 PERYLENE, BENZO(G,H,I), BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WS67 CARBAZOLE	UG/L	NA O	NA O	NA O	NA O	NA O
WV03 CHLOROMETHANE, BY GC/MS	UG/L	100 U	10 U	10 U	10 U	10 U
WV04 BROMOMETHANE, BY GC/MS	UG/L	200 U	20 U	20 U	20 U	20 U
WV05 VINYL CHLORIDE	UG/L	150 U	310 U	15 U	15 U	15 U
WV06 CHLOROETHANE, BY GC/MS	UG/L	150 U	15 U	15 U	15 U	15 U
WV07 METHYLENE CHLORIDE	UG/L	30 J	10 U	10 U	10 U	10 U
WV08 1,1-DICHLOROETHENE	UG/L	50 U	5.0 U	5.0 U	5.0 U	5.0 U
WV09 1,1-DICHLOROETHANE	UG/L	50 U	5.0 U	5.0 U	5.1 U	5.0 U

## ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 1-ADF43

VALIDATED DATA

COMPOUND	UNITS	001	002	004	005	006
WV40 TRANS-1,3-DICHLOROPROPENE	UG/L	50. U	5.00	5.00	5.00	5.00
ZZ01 SAMPLE NUMBER	NA	001	002	004	005	006
ZZ02 ACTIVITY CODE	NA	ADF43	ADF43	ADF43	ADF43	ADF43

## ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 1-ADF43

VALIDATED DATA

COMPOUND	UNITS	007	008F	009F
WD02 2,3,7,8-TCDD IN WATER (PG/ML)	PG/ML	0.000148 U		
WFO1 WATER TEMP	'C	70.8		
WFO5 PH, FIELD	SU	5.75		
WF10 CONDUCTIVITY (FIELD)	UMHOS	4150		
WH01 2,4-D	UG/L	0.2 U		
WH02 2,4,5-TP (SILVEX)	UG/L	0.2 U		
WH03 2,4,5-T	UG/L	0.2 U		
WMO1 SILVER BY ICAP	UG/L	10 U	10 U	
WMO2 ALUMINUM, TOTAL, BY ICAP	UG/L	886	75.8	
WMO3 ARSENIC, TOTAL, BY ICAP	UG/L	50 U	50 U	
WMO4 BARIUM, TOTAL, BY ICAP	UG/L	886	9.65	
WMO5 BERYLLIUM, TOTAL, BY ICAP	UG/L	2 U	2 U	
WMO6 CADMIUM, TOTAL, BY ICAP	UG/L	5 U	5 U	
WMO7 COBALT, TOTAL, BY ICAP	UG/L	10 U	10 U	
WMO8 CHROMIUM, TOTAL, BY ICAP	UG/L	12.6	10 U	
WMO9 COPPER, TOTAL, BY ICAP	UG/L	10 U	10 U	
WM10 IRON BY ICAP	UG/L	29300	50 U	
WM11 MANGANESE BY ICAP	UG/L	2860	2 U	
WM12 MOLYBDENUM BY ICAP	UG/L	NA O	NA O	
WM13 NICKEL BY ICAP	UG/L	20 U	20 U	
WM14 LEAD BY ICAP	UG/L	50 U	50 U	
WM15 ANTIMONY, TOTAL, BY ICAP	UG/L	50 U	50 U	
WM16 SELENIUM BY ICAP	UG/L	50 U	50 U	
WM17 TITANIUM BY ICAP	UG/L	NA O	NA O	
WM18 THALLIUM BY ICAP	UG/L	300 U	300 U	
WM19 VANADIUM BY ICAP	UG/L	10 U	10 U	

## ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 1-ADF43

VALIDATED DATA

COMPOUND	UNITS	007	008F	009F			
WM20 ZINC BY ICAP	UG/L	25.9	20 U				
WM21 CALCIUM, TOTAL BY ICAP	MG/L	326	2 U				
WM22 MAGNESIUM, TOTAL BY ICAP	MG/L	76.1	2 U				
WM23 SODIUM, TOTAL BY ICAP	MG/L	476	2 U				
WM24 POTASSIUM, TOTAL BY ICAP	MG/L	26.7	2 U				
WM34 MERCURY BY COLD VAPOR AA	UG/L	0.483	0.2 U				
WP01 BHC, ALPHA, BY GC/EC	UG/L	0.050 U					
WP02 BHC, BETA, BY GC/EC	UG/L	0.050 U					
WP03 BHC, DELTA	UG/L	0.050 U					
WP04 GAMMA-BHC(LINDANE), BY GC/EC	UG/L	0.050 U					
WP05 ALDRIN, BY GC/EC	UG/L	0.050 U					
WP06 DIELDRIN, BY GC/EC	UG/L	0.10 U					
WP07 ENDOSULFAN I, BY GC/EC	UG/L	0.050 U					
WP08 ENDOSULFAN II, BY GC/EC	UG/L	0.10 U					
WP09 ENDOSULFAN SULFATE, BY GC/EC	UG/L	0.10 U					
WP10 ENDRIN, BY GC/EC	UG/L	0.10 U					
WP11 ENDRIN ALDEHYDE, BY GC/EC	UG/L	NA O					
WP13 4,4'-DDE	UG/L	0.10 U					
WP14 4,4'-DDD	UG/L	0.10 U					
WP15 4,4'-DDT	UG/L	0.10 U					
WP16 TOXAPHENE	UG/L	1.0 U					
WP17 PCB-1016	UG/L	0.50 U					
WP18 PCB-1221	UG/L	0.50 U					
WP19 PCB-1232	UG/L	0.50 U					
WP20 PCB-1242	UG/L	0.50 U					
WP21 PCB-1248	UG/L	0.50 U					

## ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 1-ADF43

VALIDATED DATA

COMPOUND	UNITS	007	008F	009F		
WP22 PCB-1254	UG/L	1.0	U			
WP23 PCB-1260	UG/L	1.0	U			
WP24 CHLORDANE, TECHNICAL, BY GC/EC	UG/L	NA	O			
WP25 HEPTACHLOR, BY GC/EC	UG/L	0.050	U			
WP26 HEPTACHLOR EPOXIDE, BY GC/EC	UG/L	0.050	U			
WS01 PHENOL	UG/L	10	U			
WS03 ETHER, BIS(2-CHLOROETHYL), BY GC/MS	UG/L	10	U			
WS04 2-CHLOROPHENOL	UG/L	10	U			
WS05 1,3-DICHLOROBENZENE	UG/L	18				
WS06 1,4-DICHLOROBENZENE	UG/L	10	U			
WS07 BENZYL ALCOHOL	UG/L	10	U			
WS08 1,2-DICHLOROBENZENE	UG/L	15				
WS09 2-METHYLPHENOL (O-CRESOL)	UG/L	10	U			
WS10 ETHER, BIS(2-CHLOROISOPROPYL), BY GC/MS	UG/L	10	U			
WS11 4-METHYLPHENOL (P-CRESOL)	UG/L	10	U			
WS12 N-NITROSO-DIPROPYLAMINE	UG/L	10	U			
WS13 HEXACHLOROETHANE	UG/L	10	U			
WS14 NITROBENZENE	UG/L	10	U			
WS15 ISOPHORONE	UG/L	10	U			
WS16 2-NITROPHENOL	UG/L	10	U			
WS17 2,4-DIMETHYLPHENOL	UG/L	10	U			
WS18 BENZOIC ACID, BY GC/MS	UG/L	50	U			
WS19 METHANE, BIS(2-CHLOROETHOXY), BY GC/MS	UG/L	10	U			
WS20 2,4-DICHLOROPHENOL	UG/L	10	U			
WS21 1,2,4-TRICHLOROBENZENE	UG/L	10	U			
WS22 NAPHTHALENE	UG/L	10	U			

## ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 1-ADF43

VALIDATED DATA

COMPOUND	UNITS	007	008F	009F		
WS23 4-CHLOROANILINE	UG/L	54				
WS24 HEXACHLOROBUTADIENE,BY GC/MS	UG/L	10	U			
WS25 4-CHLORO-3-METHYLPHENOL	UG/L	10	U			
WS26 2-METHYLNAPHTHALENE	UG/L	10	U			
WS27 HEXACHLOROCYCLOPENTADIENE,BY GC/MS	UG/L	10	U			
WS28 2,4,6-TRICHLOROPHENOL	UG/L	10	U			
WS29 2,4,5-TRICHLOROPHENOL	UG/L	50	U			
WS30 2-CHLORONAPHTHALENE	UG/L	10	U			
WS31 2-NITROANILINE (ORTHO NITROANILINE)	UG/L	50	U			
WS32 PHTHALATE, DIMETHYL, BY GC/MS	UG/L	10	U			
WS33 ACENAPHTHYLENE, BY GC/MS	UG/L	10	U			
WS34 3-NITROANILINE	UG/L	50	U			
WS35 ACENAPHTHENE, BY GC/MS	UG/L	10	U			
WS36 2,4-DINITROPHENOL	UG/L	50	U			
WS37 4-NITROPHENOL	UG/L	50	U			
WS38 DIBENZOFURAN	UG/L	10	U			
WS39 2,4-DINITROTOLUENE	UG/L	10	U			
WS40 2,6-DINITROTOLUENE	UG/L	10	U			
WS41 PHTHALATE, DIETHYL, BY GC/MS	UG/L	10	U			
WS42 4-CHLOROPHENYL PHENYL ETHER	UG/L	10	U			
WS43 FLUORENE,BY GC/MS	UG/L	10	U			
WS44 4-NITROANILINE	UG/L	NA	I			
WS45 4,6-DINITRO-2-METHYLPHENOL	UG/L	50	U			
WS46 N-NITROSODIPHENYLAMINE	UG/L	10	U			
WS47 4-BROMOPHENYL PHENYL ETHER	UG/L	10	U			
WS48 HEXACHLOROBENZENE,BY GC/MS	UG/L	10	U			

## ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 1-ADF43

VALIDATED DATA

COMPOUND	UNITS	007	008F	009F
WS49 PENTACHLOROPHENOL	UG/L	50	U	
WS50 PHENANTHRENE	UG/L	10	U	
WS51 ANTHRACENE, BY GC/MS	UG/L	10	U	
WS52 PHTHALATE, DI-N-BUTYL-, BY GC/MS	UG/L	10	U	
WS53 FLUORANTHENE, BY GC/MS	UG/L	10	U	
WS54 PYRENE	UG/L	10	U	
WS55 PHTHALATE, BUTYL BENZYL	UG/L	10	U	
WS56 3,3'-DICHLOROBENZIDINE	UG/L	20	U	
WS57 ANTHRACENE, BENZO(A), BY GC/MS	UG/L	10	U	
WS58 PHTHALATE, BIS(2-ETHYLHEXYL), BY GC/MS	UG/L	19	U	
WS59 CHRYSENE, BY GC/MS	UG/L	10	U	
WS60 PHTHALATE, DI-N-OCTYL-, BY GC/MS	UG/L	10	U	
WS61 FLUORANTHENE, BENZO(B), BY GC/MS	UG/L	10	U	
WS62 FLUORANTHENE, BENZSO(K), BY GC/MS	UG/L	10	U	
WS63 PYRENE, BENZO(A), BY GC/MS	UG/L	10	U	
WS64 INDENO(1,2,3-CD)PYRENE	UG/L	10	U	
WS65 ANTHRACENE, DIBENZO(A,H), BY GC/MS	UG/L	10	U	
WS66 PERYLENE, BENZO(G,H,I), BY GC/MS	UG/L	10	U	
WS67 CARBAZOLE	UG/L	NA	O	
WV03 CHLOROMETHANE, BY GC/MS	UG/L	10.	U	10. U
WV04 BROMOMETHANE, BY GC/MS	UG/L	20.	U	20. U
WV05 VINYL CHLORIDE	UG/L	180.	U	15. U
WV06 CHLOROETHANE, BY GC/MS	UG/L	15.	U	15. U
WV07 METHYLENE CHLORIDE	UG/L	10.	U	10. U
WV08 1,1-DICHLOROETHENE	UG/L	5.0	U	5.0U
WV09 1,1-DICHLOROETHANE	UG/L	5.0	U	5.0U

## ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 1-ADF43

VALIDATED DATA

COMPOUND	UNITS	007	008F	009F		
WV10 1,2-DICHLOROETHENE, TOTAL	UG/L	500.	5.0 U	5.0U		
WV11 CHLOROFORM, BY GC/MS	UG/L	5.0 U	5.0 U	5.0U		
WV12 1,2-DICHLOROETHANE	UG/L	5.0 U	5.0 U	5.0U		
WV13 1,1,1-TRICHLOROETHANE	UG/L	5.0 U	5.0 U	5.0U		
WV14 CARBON TETRACHLORIDE, BY GC/MS	UG/L	5.0 U	5.0 U	5.0U		
WV15 BROMODICHLOROMETHANE, BY GC/MS	UG/L	5.0 U	5.0 U	5.0U		
WV16 1,2-DICHLOROPROPANE	UG/L	5.0 U	5.0 U	5.0U		
WV17 BENZENE, BY GC/MS	UG/L	130.	5.0 U	5.0U		
WV19 TRICHLOROETHENE	UG/L	5.0 U	5.0 U	5.0U		
WV20 DICHLOROPROPENE, CIS-1.3-, BY GC/MS	UG/L	5.0 U	5.0 U	5.0U		
WV21 DIBROMOCHLOROMETHANE, BY GC/MS	UG/L	5.0 U	5.0 U	5.0U		
WV22 1,1,2-TRICHLOROETHANE	UG/L	5.0 U	5.0 U	5.0U		
WV24 BROMOFORM, BY GC/MS	UG/L	5.0 U	5.0 U	5.0U		
WV25 TETRACHLOROETHENE	UG/L	5.0 U	5.0 U	5.0U		
WV26 TOLUENE	UG/L	15. J	5.0 U	5.0U		
WV27 1,1,2,2-TETRACHLOROETHANE	UG/L	5.0 U	5.0 U	5.0U		
WV28 CHLOROBENZENE, BY GC/MS	UG/L	22000. J	5.0 U	5.0U		
WV29 ETHYL BENZENE, BY GC/MS	UG/L	18. J	5.0 U	5.0U		
WV30 ACETONE, BY GC/MS	UG/L	10. U	10. U	10. U		
WV31 CARBON DISULFIDE, BY GC/MS	UG/L	5.0 U	5.0 U	5.0U		
WV32 2-BUTANONE	UG/L	10. U	10. U	10. U		
WV33 VINYL ACETATE	UG/L	10. U	10. U	10. U		
WV34 2-HEXANONE	UG/L	10. U	10. U	10. U		
WV35 4-METHYL-2-PENTANONE	UG/L	10. U	10. U	10. U		
WV36 STYRENE	UG/L	5.0U	5.0 U	5.0U		
WV37 XYLENES, TOTAL	UG/L	9.2J	5.0 U	5.0U		



## ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 1-ADF43

VALIDATED DATA

COMPOUND	UNITS	007	008F	009F		
WV40 TRANS-1,3-DICHLOROPROPENE	UG/L	5.00	5.0 U	5.00		
ZZ01 SAMPLE NUMBER	NA	007	008	009		
ZZ02 ACTIVITY CODE	NA	ADF43	ADF43	ADF43		

VALIDATED DATA

ACTIVITY ADF43      MONSANTO-QUEENY

THE PROJECT LEADER SHOULD CIRCLE ONE - STORET, AIRS, OR ARCHIVE.

CIRCLE ONE:      STORET      AIRS      ARCHIVE

FINAL DATA REPORT APPROVED BY PROJECT LEADER ON 09/18/91 17:19:19 BY Robert B. Dora.



UNITED STATES ENVIRONMENTAL PROTECTION AGENCY

REGION 7  
25 FUNSTON ROAD  
KANSAS CITY, KANSAS 66115

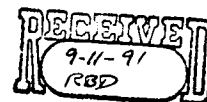
DATE: SEP 10 1991

MEMORANDUM

SUBJECT: Data Transmittal for Activity #: ADF43  
Site Description: Monsanto Queeny  
FROM: Andrea Jirka *Robert Greenall for AG*  
Chief, Laboratory Branch, ENSV  
TO: Dale Bates  
Chief, Environmental Services Division Support Branch, ENSV  
ATTN: Bob Dona

Attached is the data transmittal for the above referenced site.  
This should be considered a      Partial or X Complete data  
transmittal (completes transmittal of                                     ).  
If you have any questions or comments, please contact Dee Simmons at  
551-5129.

cc: Data Files



DRAFT

## FIELD SHEET

U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII  
 ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115

43  
 FY: P1 ACTNO: ADF13 SAMNO: 001 OCC: MEDIA: WATER PL: DONA, B.

ACTIVITY DES: MONSANTO QUEENY

LOCATION: MW-14

PROJECT NUM: A60

REF LATITUDE: -- -- --

PT: LONGITUDE: -- -- --

SAMPLE DES: Monsanto St. Louis

LOCATION: MW-14 2" PVC

CASE/PATCH/SMC: 1/1

LAB: --

DATE TIME FROM REF PT  
 BEG: 8/11/91 10:30 FAST: --  
 END: 8/11/91 : NORTH: --  
 DOWN: --

STORET/SARFAS NO: --

## ANALYSIS REQUESTED:

CONTAINER	PRESERVATIVE	MGP	NAME
CUBI	5 ML HNO3	WM	METALS
2 VOA VIALS	ICED	WV	VOLATILES
GLASS	ICED	WS	SEMIVOLATILES
GLASS	ICED	WP	PESTICIDES
GLASS	ICED	WH	HERBICIDES
CUBI	5 ML HNO3	WM34	MERCURY BY COLD VAP
GLASS	ICED	WD02	2,3,7,8-TCDD IN WATER (PGI)

COMMENTS: MW-14

Purged 11 gallons

pH-6.93

Temp-76°F

SC-1210µs/cm

Color-Black w/sheen

Strong odor

Static H<sub>2</sub>O level - 0.98 below ground surface

2" PVC well casing

PVC dedicated PVC bailer with polypropylene cord.

SAMPLE COLLECTED BY: Georgiy Miller Tom Lazarski  
Jacobs Engineering Steve Franken

DRAFT FIELD SHEET  
 U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII  
 ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115

RY: 71 ACTNO: 4071 SANNO: 002 QCC: MEDIA: WATER PL: DQNA, G.

ACTIVITY DET: KANSAS PLUMB

REF LATITUDE: -- -- --

LOCATION: TW-1

PROJECT NUM: A60

PT: LONGITUDE: -- -- --

SAMPLE DES: Menard St. Louis

DATE

TIME

FROM REF PT

LOCATION: TW-1 8" Steel

BEG: 8/18/91 17:10

EAST: -- -- --

CASE/HATCH/END: 1/1

LAR: --

END: 8/19/91

NORTH: -- -- --

STOP/START NO: --

DOWN: -- -- --

ANALYSIS REQUESTED:

CONTAINER	PRESERVATIVE	MGP	NAME
CUBI	5 ML HNO3	WM	METALS
2 VOA VIALS	ICED	WV	VOLATILES
GLASS	ICED	WS	SEMIVOLATILES
GLASS	ICED	WP	PESTICIDES
GLASS	ICED	WH	HERBICIDES
CUBI	5 ML HNO3	WM34	MERCURY BY COLD VAP
GLASS	4 C	WM02	2,3,7,8-TCDD IN WATER (PGI)

COMMENTS:

TW-1  
 Purged ~350 gallons  
 pH - 6.75  
 Temp - 75.8  
 S.C. - 2550  
 Color - Clear  
 Some color  
 Static 470 Level - 18.20 or 16.95 below ground surface  
 2" x 3' Teflon bailer with polypropylene cord  
 Casing 8" Steel w/ 570 slot S.S. Screen.

SAMPLE COLLECTED BY :

Geraghty & Miller Tom Tunnicliffe  
 Jacobs Engineering Steve Freeman

DEPT: \_\_\_\_\_ FIELD SHEET  
 U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII  
 ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115  
 FY: 91 ACTID: 4013 SAMPL: 004 VOC: \_ MEDIA: WATER PL: DONA, B.  
 ACTIVITY: MENSANTO CUSENY REF LATITUDE: \_\_\_\_\_  
 LOCATION: QS-1 PROJECT NUM: A60 PT: LONGITUDE: \_\_\_\_\_  
 SAMPLE NO: Mensanto St. Coals DATE TIME FROM REF PT  
 LOCATION: \_\_\_\_\_ BEG: 6/21/91 : : EAST: \_\_\_\_\_  
 CASE/ATCH/END: 1/1 LAB: \_\_\_\_\_ END: 6/21/91 : : NORTH: \_\_\_\_\_  
 STORET/SARDAC NO: \_\_\_\_\_ DOWN: \_\_\_\_\_

ANALYSIS REQUESTED:

CONTAINER	PRESERVATIVE	MGP	NAME
CUBI	5 ML HNO3	WM	METALS
2 VOA VIALS	ICED	WV	VOLATILES
GLASS	ICED	WS	SEMIVOLATILES
GLASS	ICED	WP	PESTICIDES
GLASS	ICED	WH	HERBICIDES
CUBI	5 ML HNO3	WM34	MERCURY BY COLD VAP
GLASS	4 C	W002	2,3,7,8-TCDF IN WATER (PG/

COMMENTS:

QS-1  
Purged - ~250 gallons one well volume 58 gallons  
PH - 8.06  
Temp - 67.1°  
S.C. - 3730  
Color - Grey  
Static H<sub>2</sub>O - 13.1 To casing lip

SAMPLE COLLECTED BY : Geraghty & Miller Tom Tannick, H  
Jacobs Engineering Steve Freeman

DRAFT FIELD SHEET  
 U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII  
 ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115

RY: 41 ACTHC: 40F16 SAMNO: 005 QCC: MEDIA: WATER PL: CCNA, B.

ACTIVITY INFO: MONSANTO CLEENY  
 LOCATION: \_\_\_\_\_

REF LATITUDE: -- -- --  
 PROJECT NUM: A60 PT: LONGITUDE: -- -- --

SAMPLE ID: 1  
 LOCATION: HW-1B  
 CASE/PATCH/ENC: 1/1  
 STREET/ROAD/ID: \_\_\_\_\_  
 DATE: 6/15/91 TIME: 14:30 FROM REF PT  
 BEG: 6/15/91 14:30 EAST: ---  
 END: 6/15/91 15:20 NORTH: ---  
 DOWN: ---

ANALYSIS REQUESTED:

CONTAINER	PRESERVATIVE	MGP	NAME
CUBI	5 HL HNO3	WM	METALS
2 VOA VIALS	ICED	WV	VOLATILES
GLASS	ICED	WS	SEMIVOLATILES
GLASS	ICED	WP	PESTICIDES
GLASS	ICED	WH	HERBICIDES
CUBI	5 HL HNO3	WM34	MERCURY BY COLD VAP
GLASS	5 C	W002	2,3,7,8-TCDD IN WATER (PG/

COMMENTS:

STATIC WATER LEVEL 18.72' BELOW TOP OF RISER  
 TOTAL DEPTH 79.67' " "  
 HEIGHT OF WATER COLUMN 60.95'  
 1 WELL VOLUME = 39.62 GALLONS

WV	WATER PH	QUALITY TEMP	PARAMS COND
0	7.88	68.7	1953
1	7.26	66.5	2020
2	6.80	65.4	2010
3	6.73	66.1	2030
4	6.60	67.4	1980

WATER IS SLIGHTLY TURBID - COLOR GRAY

SAMPLE COLLECTED BY: Klaus J. Oeder

DRAFT: FIELD SHEET  
 U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII  
 ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 65115

---

FY: 91 ACT 10: 40613 SAMPLING: 006 RCC: \_ MEDIA: WATER PL: DONA, S.

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ACTIVITY DES: NONPANTO QUEENY REF LATITUDE: \_ \_ \_  
 LOCATION: PROJECT NUM: A60 PT: LONGITUDE: \_ \_ \_

---

SAMPLE DES: VW-2B DATE TIME FROM REF PT  
 LOCATION: BEG: ~~03/11/91~~ 6/16/91 9:00 EAST: \_ \_ \_  
 CASE/LATCH/SMS: / / LAB: \_ \_ \_ END: 6/16/91 9:30 NORTH: \_ \_ \_  
 STORET/SARLAD NO: \_ \_ \_ DOWN: \_ \_ \_

ANALYSIS REQUESTED:

CONTAINER	PRESERVATIVE	MGP	NAME
CURT	5 ML HNO3	WM	METALS
2 VOA VIALS	ICED	WM	VOLATILES
GLASS	ICED	WS	SEMIVOLATILES
GLASS	ICED	WP	PESTICIDES
GLASS	ICED	WH	HERBICIDES
CURT	5 ML HNO3	WM34	MERCURY BY COLD VAP
GLASS	ICED	WM02	2,3,7,8-TCDD IN WATER (PG/

COMMENTS:

STATIC WATER LEVEL 16.49  
 1 WELL VOLUME 39 GALLONS  
 WATER QUALITY PARAS 160 GALLONS PURGED  
 WY PH TEMP COND  
 0 6.33 65.5 7761  
 1 6.12 63.7 2270  
 2 6.27 63.6 2260  
 3 6.41 63.7 2150  
 WATER SLIGHTLY TURB - COLOR GRAY

SAMPLE COLLECTED BY : Klaus J Oehler



DRAFT  
FIELD SHEET  
U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII  
ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115

RY: 41 ACTNO: 4017 SAMPLER: 007 OCC: MEDIA: WATER PL: CONA, B.

ACTIVITY DTD: MONSANTO GLENNY  
LOCATION: PROJECT NUM: 460 REF LATITUDE: PT: LONGITUDE:

SAMPLE DES: MN-8A  
LOCATION: 6/7 DATE TIME FROM REF PT  
CASE/PATCH/ONE: 1/1 LAB: REG: 03/11/91 9:15 EAST:  
STURST/SARNOG RD: END: 6/7/91 9:45 NORTH:  
DOWN:

ANALYSIS REQUESTED:  
CONTAINER PRESERVATIVE MGP NAME  
CUCI 5 ML HNO3 WM METALS  
3 VOA VIALS ICED WV VOLATILES  
GLASS ICED WS SEMIVOLATILES  
GLASS ICED WP PESTICIDES  
GLASS ICED WH HERBICIDES  
CUCI 5 ML HNO3 WM34 MERCURY BY COLD VAP  
GLASS ICED WDO2 1,2,3,7,8-TCDF IN WATER (PG/

COMMENTS:

STATIC WATER LEVEL 11.40'  
TOTAL DEPTH 48.25'  
HEIGHT WATER COLUMN 28.85  
1 WELL VOLUME 4.62  
TOTALS GALLONS PURGED 215.0  
WATER QUALITY PARAMS

WV	PH	TEMP	COND
0	5.75	70.8	4150
1	6.05	70.0	4260
2	6.51	70.1	4560
3	6.37	70.5	4670

SAMPLE COLLECTED BY: Klaus J. Cress

DRAFT FIELD SHEET  
U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII  
ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115

RY: 91 ACT NO: 4018 SAMPLING: 008 LOC: F MEDIA: WATER PL: DONA, B.

ACTIVITY DES: MONSANTO GLENNY  
LOCATION: \_\_\_\_\_

PROJECT NUM: A60 REF LATITUDE: \_\_\_\_\_  
PT: LONGITUDE: \_\_\_\_\_

SAMPLE DES: MW-9A RINS

LOCATION: \_\_\_\_\_ DATE: 6/7 TIME: FROM REF PT  
CASE/BATCH/SMC: 1/1 LAB: \_\_\_\_\_ BEG: 6/7/91 8:20 EAST: \_\_\_\_\_  
STORET/SAPDAG NO: \_\_\_\_\_ END: 6/7/91 8:20 NORTH: \_\_\_\_\_  
DOWN: \_\_\_\_\_

ANALYSIS REQUESTED:

CONTAINER	PRESERVATIVE	MGP	NAME
CURT	5 ML HNO3	WM	METALS + Hg (U.N.31)
2 VOA VIALS	ICED	WV	VOLATILES
GLASS	ICED	WS	SEMIVOLATILES
GLASS	ICED	WP	PESTICIDES
GLASS	ICED	WH	HERBICIDES
CURT	5 ML HNO3	WM34	MERCURY
GLASS	ICED	W002	2,3,7,8-TCDF IN WATER (PGI)

COMMENTS:

RINSEATE TAKEN OFF OF TEFION  
BAILER USED AT MW-9A AFTER DECON.

SAMPLE COLLECTED BY :

*Klaus J. Oeder*

CAST

## FIELD SHEET

U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII  
ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115

RY: 91 ACT93: 30P5 SAMNC: (009) OCC: F MEDIA: WATER PL: DONA, S.

ACTIVITY JES: NONNANTS CLEENY

REF LATITUDE:

LOCATION: Field Blunt

PROJECT NUM: A60

PT: LENGTH:

SAMPLE DES: 110.75970 - St Louis

LOCATION: \_\_\_\_\_

REG: ~~23121~~ 191 17:10

FROM REF PT

-----  
CASE/BATCH/NO:            /    /

445:

END:

FAST:

STORY/JARDON 17:

NORTH:

DOWN:

ANALYSIS REQUESTED:

2014-01-01

DERIVATIVE

143 P

NAME \_\_\_\_\_

2 VOW YZALS

2056

WV

2-VOLATILES

COMMENTS:

• Packaged on 06/4/91 after the sample ADF4002

SAMPLE COLLECTED BY :

DRAFT

FIELD SHEET

900P

U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII  
ENVIRONMENTAL SERVICES DIV. 225 PULSTON KANSAS CITY, KS 66115

RY: 01 ACTNO: 4010 SAMNO: ~~000~~ 000: ~~F~~ MEDIA: WATER FL: DONA, B.

ACTIVITY 100: MONSANTO QUEENY

LOCATION: USEPA (00019)

PROJECT NUM: 460

REF LATITUDE: -- --

PT: LONGITUDE: -- --

SAMPLE NO: Monsanto St. Louis

LOCATION: -- --

BEG: 03/18/91 17:30

TIME

FROM REF PT

CASE/BATCH/SNO: 1/1

LAB: --

END: 1/1

EAST: --

STORET/SAMPLE NO: --

NORTH: --

DOWN: --

ANALYST REQUESTED:

CONTAINER

PRESERVATIVE

MGP

NAME

CODI

5 mL HNO3

WN

METALS

7 VOA VIALS

1000

WV

VOLATILES

GLASS

1000

WS

SEMIVOLATILES

GLASS

1000

WP

PESTICIDES

GLASS

1000

WH

HERBICIDES

GLASS

1000

WM34

MERCURY

BY COLD VAP

GLASS

1000

WUC2

2,3,7,8-TCDF IN WATER (201)

COMMENTS:

Prepared by ESAT

Sample was in fact a PE  
Sample NOT blank!  
Analyst will report data  
Under 900P.

PE= WS 318 #10 Concent.  
the values 2 1/2 x

\*Notified C. Appleby for Organics  
6/25/91 10:50 Am.

SAMPLE COLLECTED BY: Steve Eickman

12/6/91

CONTENTS OF SHIPMENT

DESCRIPTION OF SHIPMENT	MODE OF SHIPMENT
-------------------------	------------------

PERSONNEL CUSTODY RECORD7-EPA-9262 (Revised 5/85)

✓ 6/12/91

### CONTENTS OF SHIPMENT

DESCRIPTION OF SHIPMENTMODE OF SHIPMENT

\_\_\_\_ PIECE(S) CONSISTING OF \_\_\_\_\_ BOX(ES)

COMMERCIAL CARRIER: \_\_\_\_\_

3 ICE CHEST(S); OTHER \_\_\_\_\_

~~COURIER~~  
~~SAMPLER CONVEYED~~ (SHIPPING DOCUMENT NUMBER)

## PERSONNEL CUSTODY RECORD

7-EPA-9262(Revised 5/85)

## ENVIRONMENTAL SERVICES ASSISTANCE TEAM -- ZONE II

ICF Technology Incorporated

**NSI Technology Services Inc.**

The Bionetics Corp.

ESAT Region VII  
NSI Tech.Service Inc.  
25 Funston Road  
Kansas City, KS 66115  
(913) 551-5000

---

**To:** Mike Thomas

Chief, Analytical Services Section, LABO/ENSV

**Through:** Harold Brown, Ph.D.

Deputy Project Officer for Region VII ESAT, EPA

**From:** Kevin Ludwikoski  
Scientist, ESAT



**Through:** Ron Ross

Region VII ESAT Team Leader

**Date:** 6/27/91

**Subject:** Case narrative and assignment completion.

Report:.....Monsanto-Queeny  
TID # :.....7B-9103-530  
ICF Acct. #:.....302-26-530-01  
METI Sales Order:.....1073-530  
EPA Activity #:.....ADF43  
ESAT Document Control #:.....ESAT-VII-530-0014

ESAT was assigned eight water samples for Total Metals and Mercury analysis from Monsanto-Queeny. The samples for Mercury analysis were digested and analyzed using SOP No. 3121.14 A. The Total Metals samples were digested according to SOP No. 3110.1A and analyzed on ICAP using Method 6010.

### Calibrations

All initial and continuing calibrations met Region VII guidelines for Total Metals and Mercury.

### Matrix Spike / Matrix Spike Duplicate

The recoveries for Total Metals and Mercury in the matrix spike and matrix spike duplicate met Region VII guidelines.

### Method Standard

The method standard recoveries for Total Metals and Mercury met Region VII guidelines.

### Laboratory Duplicates

The spike duplicate precision for Total Metals and Mercury met Region VII guidelines.

### Blanks

The method blank showed no Mercury above the detection limit. The method blank for Total Metals showed no analytes of interest above the detection limit with the exception of Al (73  $\mu\text{g/l}$ ). Since the concentration of Al in the samples was much higher than was found in the blank (except for sample ADF43008F), the data may be deemed acceptable for use. The samples were not blank corrected.

### PE Sample

All analytes that were present in the PE Sample were found at acceptable levels with the possible exception of Hg (69% recovery). Since no control limits had been set for this sample, and all other QC was within Region VII guidelines, this data may be deemed acceptable for use.

### Discussion

Sample ADF43003F was mistakenly labeled as a field blank when in fact it was a PE Sample prepared at the laboratory from WS 378 #10 at a 2.5X concentration. The sample has been relabeled ADF43900P and the associated true values may be found on the data sheets.

Selenium in samples ADF43001 and ADF43002 were verified by GFAA.

No problems were encountered with this project and this data may be deemed acceptable for use for its intended purpose.

### Conclusion

This assignment is now complete and data sheets for the analysis are attached. If you have any questions or comments, please contact me at 236-3881.



## ENVIRONMENTAL SERVICES ASSISTANCE TEAM -- ZONE II

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**NSI Technology Services Corporation**

The Bionetics Corp.

ESAT Region VII  
NSI Tech. Services Corp  
25 Funston Road  
Kansas City, KS 66115  
(913) 236-3881

---

TO: Robert Greenall  
ORGN Section Chief

THRU: Harold Brown, Ph.D.  
ESAT DPO

FROM: Roberta Vogel *RV*  
ESAT Scientist

THRU: Ron Ross  
ESAT Team Leader

DATE: July 3, 1991

SUBJECT: ADF43, Monsanto, Herbicides

TID.....07-9103-531  
ICF ACCOUNT #.....26-531-01  
MANTECH SALES ORDER #.....1073-531  
ESAT DOCUMENT CONTROL #...ESAT-VII-531-0021  
EPA ACTIVITY #.....ADF43

ESAT was tasked to analyze seven water samples from the Monsanto site for 2,4-D and Silvex. No target analytes were detected in these samples above the detection limit.

### HOLDING TIME

The samples were extracted by the established holding time for herbicides in water.

### INITIAL AND CONTINUING CALIBRATIONS

Five standard concentrations were used to establish linearity for 2,4-D and Silvex. Evaluation of the percent relative standard deviation of the calibration factors for these standards from their respective means revealed that both were less than 10%.

In addition two three point calibrations were run during the analysis with check calibrations using the middle level standard midway between calibrations. The % differences for all of the check calibrations were less than 25%.

## **QUANTITATION**

Peak heights from the DB5 column were used for quantitation. Quantitation was performed using the mean calibration factor from the nearest calibration for each analyte.

## **METHOD BLANK**

No target analytes were detected in the method blank.

## **QC SPIKES**

A method standard and performance evaluation sample were analyzed. In the method standard 2,4-D and Silvex were recovered at 123 and 83% respectively. In the performance evaluation sample 2,4-D and Silvex were recovered at 85 and 62% respectively. There are no established guidelines for recovery of herbicide spikes.

## **RESULTS AND DISCUSSION**

One confirmed positive hit for Silvex was found in sample 004. However, the quantitations differed by 1000% between columns, so the result was reported as a non-detect at the lower level detected, as directed in the organic analysis SOP. The data has been entered into last.

## ENVIRONMENTAL SERVICES ASSISTANCE TEAM -- ZONE II

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Kansas City, KS 66115  
(913) 236-3881

---

TO: Robert Greenall  
ORGN Section Chief

THRU: Harold Brown, Ph.D.  
ESAT DPO

FROM: Roberta Vogel *Ø*  
ESAT Scientist

THRU: Ron Ross  
ESAT Team Leader

DATE: July 8, 1991

SUBJECT: ADF43, Monsanto, Herbicides

TID.....07-9103-531  
ICF ACCOUNT #.....26-531-01  
MANTECH SALES ORDER #.....1073-531  
ESAT DOCUMENT CONTROL #...ESAT-VII-531-0021  
EPA ACTIVITY #.....ADF43

This is an amendment to the original ADF43 report dated 7/3/91. In the original report, data regarding the analysis for 2,4,5-T was not included. It has been added to the LAST report and to the data package. 2,4,5-T was not detected in any of the samples and was recovered at 96% in the matrix spike. The LAST data is attached.

## ENVIRONMENTAL SERVICES ASSISTANCE TEAM -- ZONE II

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NSI Technology Services  
25 Funston Road  
Kansas City, KS 66115  
(913) 236-3881

---

To: Robert Greenall  
Chief, ORGN, LABO, EPA Region VII

From: Janet K. Muse *JM*  
Senior Scientist, ESAT

Thru: Ronald A. Ross *RAM for RAR*  
Region VII ESAT Team Leader

Thru: Harold Brown, Ph.D.  
Deputy Project Officer, Region VII ESAT, EPA

Date: July 3, 1991

Subject: Case Narrative and Data Review Document, **Monsanto**

TID #: ..... 07-9103-531  
ICF Acct. #: ..... 302-26-531-01  
NSI Sales Order : ..... 1073-531  
EPA Activity #: ..... **ADF43**  
ESAT Document Control #: ..... ESAT-VII-531-0018

Seven water samples were received by ESAT for BNA analysis.

### GC/MS Performance

All Region VII GC/MS performance criteria were met for a 50 nanogram injection of DFTPP.

### Method Blanks

Bis(2-ethylhexyl)phthalate was present in the method blank at 3.1 ug/L. Positive hits for this compound were only reported positive in the samples after the blank rule was applied. That is, positive hits for bis(2-ethylhexyl)phthalate in the samples were reported and coded with a "U" unless they were greater than 10-times the level in the method blank.

### Initial Calibration

All compounds met the 0.05 criteria for response factors. All compounds were less than 30% for RSD in the initial calibration.

### **Continuing Calibration**

4-nitroaniline was less than the 0.05 criteria limit for response factors in the continuing calibration of 6/25/91. 4-nitroaniline was invalidated in all the samples analyzed on that day.

A few compounds were outside the 35% difference for response factors in the continuing calibrations. The only positive hit requiring data qualification was 4,6-dinitro-2-methylphenol in ADF43003F. 4,6-dinitro-2-methylphenol was J-coded due to this calibration outlier.

### **Surrogate Recoveries**

Recoveries for surrogates were within the established Region VII guidelines, except for 2-fluorobiphenyl in all the samples. 2-fluorophenol was the only acid surrogate outside criteria limits in ADF43002. No data was qualified. Most of the surrogates were diluted out of ADF43001, -001S, and -001W.

### **Matrix Spike / Matrix Spike Duplicate**

A matrix spike and matrix spike duplicate were analyzed for ADF43001. Due to matrix effects in sample ADF43001, the MS/MSD had to be diluted 1:10. Because of this, most of the spiked compounds were diluted out. No data was qualified.

### **Summary**

From the evaluation of the QC mentioned above, this data meets the requirements for acceptability.

This analysis is now complete. Data sheets are attached. If you have any questions concerning this analysis, contact me.

**ICF Technology Incorporated  
NSI Technology Services Corporation**

**EPA Region VII Laboratory  
NSI Tech. Services Corp.  
25 Funston Rd.  
Kansas City, KS 66115  
(913) 236-3881**

**To:** Robert Greenall  
Chief, ORGN, LABO, Region VII EPA

**From:** Laleh A. Jonooby *AJ*  
Senior Scientist, ESAT

Paula Woodland *AJ for PW*  
Chemist, ESAT

**Thru:** Ronald A. Ross  
Region VII ESAT Team Leader

**Thru:** Harold Brown, Ph.D.  
Deputy Project Officer, Region VII ESAT, EPA

**Date:** June 27, 1991

**Subject:** Case Narrative and Data Review Document, **MONSANTO**

TID #: ..... 7B-9103-531  
ICF Acct. #: ..... 302-26-531-01  
NSI Sales Order : ..... 1073-531  
EPA Activity #: ..... **ADF43**  
ESAT Document Control #:... ESAT-VII-531-0017

Nine water samples were received by ESAT for the analysis of volatile organic compounds.

Samples arrived in the lab with headspace.

**Holding Time**

The 14 day holding time for all volatile compounds were met for all of the water samples except for the dilution of sample ADF43-007 (1:100). Chorobenzene in that sample was J-coded.

**GC/MS performance**

All Region VII GC/MS performance criteria were met for a 50 nanogram injection of 4-bromofluorobenzene.

## **Method Blank**

A volatile method blank was analyzed each day samples were run. Positive hits for target compounds in the blanks were only reported positive in the sample after the blank rule was applied. That is, positive hits for 2-butanone in the samples were reported and coded with a "U" unless they were greater than 10-times the level in the daily blank. A 5-times rule was applied for the other analytes.

## **Initial Calibration**

Two initial calibrations, 20 - 200 ug/L were analyzed for this activity. The relative percent difference limit of 30% for response factors was met for all analytes in the initial calibration of 6/11/91 with the exception of methylene chloride which was outside of the 30% criteria limit. Methylene chloride in sample -001 was J-coded. Only four analytes were outside the criteria limits in the second initial calibration. No data was qualified by the second initial calibration.

## **Continuing Calibration**

The criteria limits for response factor of 0.30 and the percent difference of 35% between response factors were met for all analytes present in the samples in all of the continuing calibrations with the exception of methylene chloride being outside the criteria limits in the continuing calibration of 6/13. However, methylene chloride was already J-coded due to the initial calibration.

## **Surrogate Recovery**

Surrogate recoveries in the samples and blanks were within the established control limits for this analysis with the exception of sample -007 in which all three surrogate recoveries were outside the criteria limits. Toluene, ethylbenzene and total xylenes were J-coded in sample -007.

## **Matrix Spike/Matrix Spike Duplicate**

A matrix spike and a matrix spike duplicate were analyzed for sample ADF43-004. Spike recoveries were within normal limits for all of the analytes except for 4-methyl-2-pentanone and 2-hexanone which had high recoveries in both of the MS and MSD. This is probably due to contamination within the purge and trap system from previous analyses since neither of the two compounds were present in any of the samples. No data was qualified by the MS/MSD.

## **Discussion**

Samples arrived in the lab with headspace. The affect of this would be to bias all data low and non-detect data is suspect.

Chloroform was J-coded in ADF43-003F, due to the quantitation being outside the calibration range. Small quantities of chlorobenzene were found in sample -006 below the method detection limits.

003F was  
marked wrong  
actually J 960  
(PESam  
dx

Library searches of all of the water samples were performed. See the attached data sheets.

### **Summary**

From the evaluation of the QC mentioned above, this data meets the requirements for acceptability.

This analysis is now complete with data sheets attached. If you have any questions please feel free to contact us.



## ENVIRONMENTAL SERVICES ASSISTANCE TEAM -- ZONE II

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(913) 551-5000

---

TO: Robert Greenall  
Chief, ORGN, LABO/ENSV

FROM: Louis Tollackson *LT*  
ESAT

THRU: Ron Ross  
Region VII ESAT Team Leader, NSI-TSC-ES

THRU: Harold Brown, Ph.D.  
Deputy Project Officer for Region VII ESAT, EPA

DATE: July 29, 1991

SUBJECT: Case Narrative - Monsanto

TID # ..... 07-9103-531  
ICF Acct. #: ..... 26-531-01  
NSI #: ..... 1073-531  
EPA Activity #: ..... ADF43  
ESAT Document Control #: ..... ESAT-VII-531-0023

Six water samples, ADF43001-002, and ADF43004-007 were recieved by ESAT on 6/12/91 for pesticide analysis. The samples were extracted on 6/17/91. The sample extracts were initially analyzed by GLC on 7/7/91. Evaluation of the data determined that the analysis did not meet QC criteria due to column and detector contamination from a previous activity. The instrument was cleaned, and the sample extracts were reanalyzed on 7/10-11/91. The analysis results gave an unconfirmed indication for aldrin which may have been due to dimethlypthalate contamination. Because of low DCB recoveries, samples ADF30002, ADF30004, and ADF30005 were reextracted on 7/18/91 and analyzed by GLC on 7/22/91. The results were negative for pesticides and PCB's.

### **Holding Time**

The water sample was extracted within the seven day holding time established for water.

### **Calibration and Continuing Calibration**

Quantitation was done by a three point calibration and results were found by linear regression analysis adjusted for dilution, final volume, amount of sample injected, and volume of sample extracted. Continuing calibration was checked by analyzing a mid level standard every five sample injections. The D% range for these standards was from 1% - 20% for all analytes of interest. The continuing calibration meets Region VII Guidelines. Continuing calibration for the reextracts was 1 - 9 %D which is within Region VII Guidelines for continuing calibrations.

### **Method Blank**

A method blank was analyzed and contained no analytes of interest for this analysis.

### **Method Standard**

A method standard was analyzed for this analysis. G-BHC: 95% recovery, Heptachlor: 58% recovery, Dieldrin: 82% recovery, Endrin: 90% recovery, and DDT: 90% recovery from the matrix spike. These values are within Region VII Guidelines for matrix spike recoveries for water. Due to the possibility of dimethylphthalate contamination of either reagent or glassware, the aldrin recovery could not be calculated. Method standard data is not used to qualify sample results. A method standard was also analyzed for the reextracts and yielded the following results: G-BHC: 56% recovery, Heptachlor: 56% recovery, Dieldrin: 71% recovery, Endrin: 65%. Aldrin was not calculated due to the presence of an interferent.

### **Performance Standard**

A performance standard was analyzed. The recoveries for the extraction were: Heptachlor - 117%, Aldrin - 71%, Dieldrin - 132%, DDD - 100%. These values are within Region VII Guidelines for matrix spike recoveries.

### **Matrix Spike**

A matrix spike was analyzed with the reextracted water samples. The recovery results are G-BHC - 66%, Heptachlor - 80%, Dieldrin - 70% and Endrin - 74%.

### **Surrogate Recoveries**

DCB (Decachlorobiphenyl) was the surrogate added to the sample

extracts. The DCB recoveries varied with this analysis. The recoveries ranged from 14% to 104%. Samples ADF43002, ADF43004, and ADF43005 yielded 13% to 19% DCB recoveries. The recoveries for all other samples were within Region VII Guidelines. Explanations for these low recoveries including leaky equipment, or overheating of sample extract during sample concentration and solvent exchange. Samples ADF43002, ADF43004, and ADF439005 were reextracted on 7/18/91. The range of DCB recoveries for these reextracts were 94% - 111%. The recoveries are within Region VII Guidelines for DCB recoveries.

## **Results and Discussion**

Six water samples were analyzed for pesticides. The combined results were negative for pesticides (See data sheets). Three samples exhibited DCB recoveries lower than normal. They were reextracted and the analysis results were negative. The samples also contained an unconfirmed aldrin like compound which may be dimethylphthalate a reagent or glassware contaminate. The analysis is now complete. If you have any questions, please contact ESAT.

## ENVIRONMENTAL SERVICES ASSISTANCE TEAM -- ZONE II

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---

To: Robert Greenall  
Chief, ORGN, LABO, EPA Region VII

From: Kirit H Bhatt *KHB*  
Senior Organic Analytical Chemist, ESAT

Thru: Ronald A. Ross  
Region VII ESAT Team Leader

Thru: Harold Brown, Ph.D.  
Deputy Project Officer, Region VII ESAT, EPA

Date: July 30, 1991

Subject: Case Narrative and Data Review Document, Monsanto

TID #: ..... 07-9103-531  
ICF Acct. #: ..... 302-26-531-01  
NSI Sales Order : ..... 1073-531  
EPA Activity #: ..... ADF43  
ESAT Document Control #: ..... ESAT-VII-531-0027

ESAT was assigned to extract and analyse six water samples and one field blank for 2,3,7,8-TCDD, by HRGC-HRMS. The analysis of matrix spikes ADF43900G and ADF43901G could not be concluded due to an electronic fault of the HRMS.

### Holding Times

The seven day holding time from sampling to extraction and the 40 day holding time from extraction to analysis were met for all water samples.

### GC/MS Performance

HRGC-HRMS performance was accertained by analysing a continuous calibration standard, which met the specified response factor requirement with respect to the initial calibration.

### Method Blanks

The method blank ADF43900M was analysed twice during the analysis of the activity. Both analysis showed prescence of low concentration of 2,3,7,8-TCDD. Blank solvent injections were carried out before analysing the method and field blanks. No peaks were detected. Consequently, the blank rule(5xsignal for the analyte)was applied. The final tabulated results are attached.

**Initial Calibration**

Initial calibration of 2,3,7,8-TCDD was linear, with mean RF of 1.29.

**Continuous Calibration**

Continuous calibration was within the range of the initial calibration, (RF=1.28).

**Surrogate Recoveries**

Recoveries for surrogates ranged between 104% - 116%. No limits for these recoveries are specified in the EPA protocol.

**Matrix Spike/Duplicate Spike**

Matrix spikes were extracted but not analysed due to the electronic fault of the HRMS.

**Performance Evaluation**

No performance evaluation sample was analyzed with this activity.

**Summary**

From the evaluation of the QC mentioned above, this data meets the requirements for acceptability.

This analysis is now complete. Please contact me if there are any queries about this activity.

The HRGC-HRMS analysis was jointly carried out by myself and Shaher Mohshin(EPA).

5m

QUALITY CONTROL EXCEPTIONS/EXPLANATIONS

1) The following samples have been J-Coded for the reasons given:

<u>Sample No.</u>	<u>Parameter(s)</u>	<u>Reason</u>
ADF43-007	Toluene, ethylbenzene, <del>styrene</del>	Surrogates were outside criteria limits
-007	chlorobenzene	H.T. time passed
-001	CH <sub>2</sub> Cl <sub>2</sub>	cut on cent. cal
-003F *	Chloroform	Outside calibration range

2) The following quality control exceptions are included in this data package.

\* 003F sample marked in error  
actually P.E. sample (900P) ds

# TENTATIVELY IDENTIFIED COMPOUNDS

TITLE: MONSANTO

LAB:

ANALYST/ENTRY: LAJ

REVIEW LEVEL: 2

MATRIX: WATER

METHOD: 5241D00

REVIEWER: *Aj for Jm*

DATA FILE: *Aj* LJ2

UNITS: UG/L

CASE:

DATE: 06/27/93

SAMPLE NO.	COMPOUND NAME	FRACTION	EST. CONCENTRATION
ADF43007			
	ALKANE	VOA	6 J
	SUBSTITUTED BENZENE	VOA	12 J
ADF43008F			
	UNKNOWN	VOA	85 J

- \* THIS IS A CRUDE ESTIMATION BASED ON RESPONSE RELATIVE TO AN INTERNAL STANDARD. AN AUTHENTIC STANDARD HAS NOT BEEN RUN.
- \*\* THE COMPOUNDS WERE IDENTIFIED USING A LIBRARY SEARCH ROUTINE. AUTHENTIC STANDARDS HAVE NOT BEEN ANALYZED TO VERIFY COMPOUND MASS SPECTRA AND RETENTION TIMES.